

Full wwPDB X-ray Structure Validation Report (i)

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PDB ID : 1GAM

Title : GAMMA B CRYSTALLIN TRUNCATED C-TERMINAL DOMAIN

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Deposited on : 1996-02-02

Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org A user guide is available at

https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity: 4.02b-467

Xtriage (Phenix) : NOT EXECUTED EDS : NOT EXECUTED

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

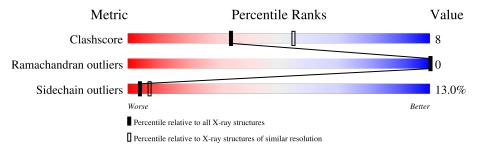
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain		
1	A	86	76%	20%	5%
1	В	86	74%	20%	5% •



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1489 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

• Molecule 1 is a protein called GAMMA B CRYSTALLIN.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	86	Total 726	C 455	N 134	O 131	S 6	0	0	0
			Total			0	$\frac{\sigma}{S}$			
1	1 B	86	726		134	131	6	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	ALA	ASP	conflict	UNP P02526
В	107	ALA	ASP	conflict	UNP P02526

• Molecule 2 is water.

Mo	ol	Chain	Residues	Atoms	ZeroOcc	AltConf
2		A	23	Total O 23 23	0	0
2		В	14	Total O 14 14	0	0

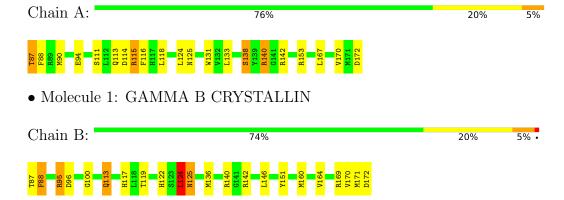


3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

• Molecule 1: GAMMA B CRYSTALLIN





4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source	
Space group	P 32 2 1	Depositor	
Cell constants	36.48Å 36.48Å 232.80Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor	
Resolution (Å)	8.00 - 2.60	Depositor	
% Data completeness	(Not available) (8.00-2.60)	Depositor	
(in resolution range)	(1101 available) (0.00 2.00)		
R_{merge}	0.07	Depositor	
R_{sym}	(Not available)	Depositor	
Refinement program	X-PLOR	Depositor	
R, R_{free}	0.210 , (Not available)	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	1489	wwPDB-VP	
Average B, all atoms (Å ²)	17.0	wwPDB-VP	



5 Model quality (i)

5.1 Standard geometry (i)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with |Z| > 5 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.55	0/743	0.82	0/998	
1	В	0.57	0/743	0.88	2/998 (0.2%)	
All	All	0.56	0/1486	0.85	2/1996 (0.1%)	

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\mathrm{Ideal}(^{o})$
1	В	124	LEU	CA-CB-CG	9.33	136.75	115.30
1	В	88	PHE	N-CA-C	-5.44	96.31	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	726	0	700	14	0
1	В	726	0	700	13	1
2	A	23	0	0	1	0
2	В	14	0	0	1	2
All	All	1489	0	1400	24	3

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (24) close contacts within the same asymmetric unit are listed below, sorted by their clash



magnitude.

A.1 1	A. 0	Interatomic	Clash
Atom-1	Atom-2	$\operatorname{distance}\ (\text{\AA})$	overlap (Å)
1:B:95:ARG:HG2	1:B:95:ARG:HH11	1.35	0.91
1:A:142:ARG:HD3	1:B:142:ARG:HD2	1.71	0.71
1:B:136:MET:HE1	1:B:140:ARG:HH21	1.56	0.71
1:B:95:ARG:HG2	1:B:95:ARG:NH1	2.07	0.62
1:A:142:ARG:CD	1:B:142:ARG:HD2	2.31	0.58
1:A:87:THR:HG23	1:A:87:THR:O	2.03	0.58
1:B:87:THR:O	1:B:87:THR:HG22	2.04	0.57
1:A:94:GLU:HG3	1:A:118:LEU:CD2	2.35	0.56
1:B:124:LEU:HD22	1:B:151:TYR:HB2	1.92	0.51
1:A:170:VAL:HG11	1:B:170:VAL:HG11	1.95	0.48
1:B:88:PHE:CG	1:B:169:ARG:HD2	2.49	0.48
1:A:115:ARG:HG2	1:A:116:PHE:CE2	2.51	0.46
1:A:88:PHE:HB3	1:A:131:TRP:CZ2	2.52	0.45
1:A:90:MET:CE	1:A:124:LEU:HD13	2.47	0.45
1:A:94:GLU:HG3	1:A:118:LEU:HD21	1.97	0.45
1:A:87:THR:O	1:A:87:THR:CG2	2.65	0.45
1:A:133:LEU:HA	1:A:167:LEU:HD23	2.00	0.44
1:A:138:SER:O	1:A:140:ARG:HG2	2.20	0.42
1:B:96:ASP:OD1	1:B:122:HIS:HB3	2.20	0.42
1:B:95:ARG:HG3	2:B:186:HOH:O	2.20	0.41
1:B:125:ASN:HD22	1:B:125:ASN:C	2.23	0.41
1:B:113:GLN:O	1:B:117:HIS:HA	2.20	0.41
1:A:87:THR:N	2:A:189:HOH:O	2.54	0.40
1:A:90:MET:HE3	1:A:124:LEU:HD13	2.02	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
2:B:176:HOH:O	2:B:176:HOH:O[6_765]	1.93	0.27
2:B:185:HOH:O	2:B:185:HOH:O[6_765]	1.95	0.25
1:B:100:GLY:O	1:B:117:HIS:ND1[6_765]	2.12	0.08

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries



of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
1	A	84/86 (98%)	82 (98%)	2 (2%)	0	100	100
1	В	84/86 (98%)	82 (98%)	2 (2%)	0	100	100
All	All	168/172 (98%)	164 (98%)	4 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	77/77 (100%)	67 (87%)	10 (13%)	4 7
1	В	77/77 (100%)	67 (87%)	10 (13%)	4 7
All	All	154/154 (100%)	134 (87%)	20 (13%)	4 7

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	THR
1	A	111	SER
1	A	113	GLN
1	A	114	ASP
1	A	115	ARG
1	A	125	ASN
1	A	138	SER
1	A	140	ARG
1	A	153	ARG
1	A	172	ASP
1	В	95	ARG
1	В	113	GLN
1	В	119	THR
1	В	124	LEU

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Mol	Chain	Res	Type
1	В	125	ASN
1	В	146	LEU
1	В	160	MET
1	В	164	VAL
1	В	171	MET
1	В	172	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	ASN
1	В	125	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

